GAUGE-FIELD THEORIES

Ulrich Mosel
Institut für Theoretische Physik, Universität Giessen
D–35392 Giessen, Germany

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Contents

Part I

Preliminaries

Chapter 1

UNITS AND METRIC

1.1 Units

It is customary in elementary particle physics and field theories to choose the system of units such that the resulting expressions assume a simple form. In particular, instead of working with the usual three mechanical units for length, mass and time one introduces instead three basic units for velocity, action and length.

The choice of the velocity of light, c, as the unit for the velocity implies that

$$c = 1. \tag{1.1}$$

This in turn means that in such a system length and time have the same dimensions and are equivalent units. With $c = 1$ the relativistic energy-momentum relation assumes the simple form

$$E^2 = p^2 + m^2. \tag{1.2}$$

Choosing next the unit of action, $\hbar$, such that

$$\hbar = 1 \tag{1.3}$$

connects the dimensions of mass, time and length. Since time and length are equivalent mass assumes the dimension of an inverse length; the same holds consequently for the momentum and the energy (equation 1.2). The choice of 1 fm $= 10^{-13}$ cm as a length unit, for example, leads to masses, momenta and energies all given in fm^{-1}. The transformation of the expressions back to the standard MKSA system can be achieved by multiplying all quantities with the appropriate combinations of $\hbar$ and c.

1.2 Metric and Notation

Spatial three-vectors in this book are written with an overhead arrow; in contrast, vectors in some internal space are distinguished by boldface print.

In relativistic expressions four-vectors are always written as

$$A^\mu = (A^0, \vec{A}) = (A^0, A^1, A^2, A^3) \quad . \tag{1.4}$$

Four-vectors with an upper index are called *contravariant* vectors. *Covariant* vectors, denoted by lower indices, are then defined by

$$A_\mu = g_{\mu\nu} A^\nu \tag{1.5}$$

with the *metric tensor*

$$g_{\mu\nu} = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & -1 & 0 & 0 \\ 0 & 0 & -1 & 0 \\ 0 & 0 & 0 & -1 \end{pmatrix} \quad . \tag{1.6}$$

The metric tensor with upper indices is defined as the inverse of g. It is given by

$$g^{\mu\nu} = g_{\mu\nu} \ . \tag{1.7}$$

A scalar product of 2 four-vectors is then defined by

$$A \cdot B = A_\mu B^\mu = A^\mu B_\mu \ . \tag{1.8}$$

The invariant square of the four-vector is thus given by

$$A^2 = A_\mu A^\mu = A_0^2 - \vec{A}^2 \quad . \tag{1.9}$$

Here, as always in the text, the Einstein convention of summing over double indices, one lower and one upper, is used. The components of four-vectors are generally labeled by Greek indices, whereas roman indices are used to refer specifically to the last three (vector) components.

The space-time four-vector is

$$x^\mu = (x^0, \vec{x}) = (t, \vec{x}) \ , \tag{1.10}$$

and the four-momentum is given by

$$p^\mu = (E, \vec{p}) \ . \tag{1.11}$$

Finally, the components of the contravariant four-gradient are abbreviated by

$$\partial^\mu \equiv \frac{\partial}{\partial x_\mu} = \left(\frac{\partial}{\partial t}, -\vec{\nabla}\right) \ , \tag{1.12}$$

so that the four-momentum operator is given by

$$\hat{p}^\mu = \mathrm{i}\partial^\mu \ , \tag{1.13}$$

and the Lorentz-invariant *d'Alembert operator* by

$$\partial^\mu \partial_\mu = \frac{\partial^2}{\partial t^2} - \vec{\nabla}^2 = \Box \ . \tag{1.14}$$

Other important four-vectors are the current

$$j^\mu = \left(\rho, \vec{j}\right) \tag{1.15}$$

and the electromagnetic potential

$$A^\mu = \left(\phi, \vec{A}\right) \ . \tag{1.16}$$

Part II

Fundamentals of Field Theory

Chapter 2

CLASSICAL FIELDS

In this chapter a few essential facts of classical field theory will be discussed. It will first be shown how to derive the equations of motion of a field theory, for example the Maxwell equations of electrodynamics, from a Lagrangian. Secondly, the connection between symmetries of the Lagrangian and conservation laws will be discussed. This is one of the central points for all the theoretical developments in this book.

2.1 Equations of Motion

The equations of motion of classical mechanics can be obtained from a Lagrange function by using Hamilton's principle that the action for a given mechanical system is minimal for the physical space-time development of the system.

The equations of motion for fields that determine their space-time dependence can be obtained in an analogous way by identifying the field amplitudes at a coordinate $\vec{x}$ with the dynamical variables (coordinates) of the theory.

Let the functions that describe a field be denoted by

$$\Phi_\alpha(x) \quad \text{with} \quad x^\mu = (t, \vec{x}) \ . \tag{2.1}$$

The Lagrangian L of the system is expressed in terms of a Lagrange density $\mathcal{L}$, as follows:

$$L = \int \mathcal{L}(\Phi_\alpha, \partial_\mu \Phi_\alpha) d^3x \tag{2.2}$$

where the spatial integration is performed over the volume of the system.

The action S is then defined as usual by

$$S = \int_{t_0}^{t_1} L \; dt = \int_{\Omega} \mathcal{L}(\Phi_\alpha, \partial_\mu \Phi_\alpha) d^4x \tag{2.3}$$

with $d^4x = d^3x \; dt$. The, in general finite, space-time volume of the system is denoted by Ω.

As pointed out before the fields $\Phi_\alpha(x)$ play the same role as the generalized coordinates q_i in classical mechanics; the analogy here is such that the fields Φ_α correspond to the coordinates q and the points $\vec{x}$ to the classical indices i. The corresponding velocities are in a direct analogy given by the time derivative of $\Phi_\alpha : \partial_t \Phi_\alpha$. Lorentz covariance then requires that also the derivatives with respect to the first three coordinates appear; this explains the presence of the four-gradients $\partial_\mu \Phi_\alpha$ in eq. (2.3).

The action S is now varied with respect to the fields and their derivatives:

$$\begin{aligned} \Phi_\alpha(x) &\rightarrow \Phi'_\alpha(x) = \Phi_\alpha(x) + \delta\Phi_\alpha(x) \\ \partial_\mu \Phi_\alpha(x) &\rightarrow (\partial_\mu \Phi_\alpha(x))' = \partial_\mu \Phi_\alpha(x) + \delta(\partial_\mu \Phi_\alpha(x)) \quad . \end{aligned} \tag{2.4}$$

This yields

$$\begin{aligned} \delta S &= \int_{\Omega} \mathcal{L}(\Phi'_\alpha, (\partial_\mu \Phi_\alpha)') d^4x - \int_{\Omega} \mathcal{L}(\Phi_\alpha, \partial_\mu \Phi_\alpha) d^4x \\ &= \int \left(\frac{\partial \mathcal{L}}{\partial \Phi_\alpha} \delta\Phi_\alpha + \frac{\partial \mathcal{L}}{\partial(\partial_\mu \Phi_\alpha)} \delta(\partial_\mu \Phi_\alpha) \right) d^4x \\ &= \int \left[\left(\frac{\partial \mathcal{L}}{\partial \Phi_\alpha} - \partial_\mu \frac{\partial \mathcal{L}}{\partial(\partial_\mu \Phi_\alpha)} \right) \delta\Phi_\alpha \right. \\ &\qquad \left. + \partial_\mu \left(\frac{\partial \mathcal{L}}{\partial(\partial_\mu \Phi_\alpha)} \delta\Phi_\alpha \right) \right] d^4x \quad . \end{aligned} \tag{2.5}$$

According to the Einstein-convention a summation over μ is implicitly contained in this expression.

The equations of motion are now obtained from the variational principle

$$\delta S = 0 \tag{2.6}$$

for arbitrary variations $\delta\Phi_\alpha$ under the constraint that

$$\delta\Phi_\alpha(t_0) = \delta\Phi_\alpha(t_1) = 0 \quad .$$

The last term in eq. (2.5) can be converted into a surface integral by using Gauss's law; for fields which are localized in space this surface integral vanishes. Since the variations $\delta\Phi_\alpha$ are arbitrary the condition $\delta S = 0$ leads to the equations of motion

$$\frac{\partial}{\partial x^\mu}\left(\frac{\partial \mathcal{L}}{\partial(\partial_\mu \Phi_\alpha)}\right) - \frac{\partial \mathcal{L}}{\partial \Phi_\alpha} = 0 \quad . \tag{2.7}$$

The relativistic equivalence principle demands that these equations have the same form in every inertial frame of reference, i.e. that they are Lorentz-covariant. This is only possible if $\mathcal{L}$ is a Lorentz scalar, i.e. if it has the same functional dependence on the fields and their derivatives in each reference frame.

In a further analogy to classical mechanics, the canonical field momentum is defined as

$$\Pi_\alpha = \frac{\partial \mathcal{L}}{\partial \dot{\Phi}_\alpha} = \frac{\partial \mathcal{L}}{\partial(\partial_0 \Phi_\alpha)} \quad . \tag{2.8}$$

From $\mathcal{L}$ and Π_α the Hamiltonian H is obtained as

$$H = \int \mathcal{H}\, d^3x = \int (\Pi_\alpha \dot{\Phi}_\alpha - \mathcal{L}) d^3x \quad . \tag{2.9}$$

The Hamiltonian H represents the energy of the field configuration.

2.1.1 Examples

2.1.1.1 Electrodynamics

The two inhomogeneous Maxwell equations[1]

$$\vec{\nabla} \cdot \vec{E} = \rho \tag{2.10}$$

$$\vec{\nabla} \times \vec{B} - \frac{\partial \vec{E}}{\partial t} = \vec{j} \tag{2.11}$$

can be rewritten by introducing the four-vector field

$$A^\mu = (A^0, \vec{A}) \ , \tag{2.12}$$

defined via

$$\begin{aligned} \vec{B} &= \vec{\nabla} \times \vec{A} \\ \vec{E} &= -\vec{\nabla} A^0 - \frac{\partial \vec{A}}{\partial t} \quad . \end{aligned} \tag{2.13}$$

[1] Here the Heaviside units are used with $c = 1$ and $\epsilon_0 = \mu_0 = 1$.

Note that the 2 homogeneous Maxwell equations are now automatically fulfilled.

In terms of A^μ the antisymmetric field tensor

$$F^{\mu\nu} = \frac{\partial A^\nu}{\partial x_\mu} - \frac{\partial A^\mu}{\partial x_\nu} = \partial^\mu A^\nu - \partial^\nu A^\mu \tag{2.14}$$

can be defined. With its help the equations of motion (2.10,2.11) can be written as

$$\partial_\mu F^{\mu\nu} = \frac{\partial F^{\mu\nu}}{\partial x^\mu} = j^\nu \quad , \tag{2.15}$$

with the four-current

$$j^\nu = (\rho, \vec{j}) \quad . \tag{2.16}$$

Equation (2.15) is the equation of motion for the field tensor or the four-vector field A_μ.

It is easy to show that eq. (2.15) can be obtained from the Lagrangian

$$\mathcal{L} = -\frac{1}{4} F_{\mu\nu} F^{\mu\nu} - j^\nu A_\nu \tag{2.17}$$

by using equation (2.7); the fields A^ν here play the role of the fields Φ_α in (2.7). We have

$$\begin{aligned} \frac{\partial \mathcal{L}}{\partial A_\nu} &= -j^\nu \\ \frac{\partial \mathcal{L}}{\partial(\partial_\mu A_\nu)} &= -\frac{1}{4} 2(+F^{\mu\nu} - F^{\nu\mu}) = -F^{\mu\nu} \quad . \end{aligned} \tag{2.18}$$

The last step is possible because F is an antisymmetric tensor.

The equation of motion is therefore

$$\frac{\partial}{\partial x^\mu}\left(\frac{\partial \mathcal{L}}{\partial(\partial_\mu A_\nu)}\right) - \frac{\partial \mathcal{L}}{\partial A_\nu} = -\frac{\partial F^{\mu\nu}}{\partial x^\mu} + j^\nu = 0 \quad , \tag{2.19}$$

in agreement with eq. (2.15). It is now easy to interpret the two terms in $\mathcal{L}$ (eq. (2.17)): the first one gives the Lagrangian for the free electromagnetic field, whereas the second describes the interaction of the field with charges and currents.

The two homogeneous Maxwell equations can also be expressed in terms of the field tensor by first introducing the *dual field tensor*

$$\tilde{F}^{\mu\nu} = \frac{1}{2} \varepsilon^{\mu\nu\rho\sigma} F_{\rho\sigma} \; ; \tag{2.20}$$

here $\varepsilon^{\mu\nu\rho\sigma}$ is the Levi-Civita antisymmetric tensor which assumes the values +1 or -1 according to whether $(\mu\nu\rho\sigma)$ is an even or odd permutation of (0,1,2,3), and 0 otherwise. In terms of $\tilde{F}$ the homogeneous Maxwell equations read

$$\partial_\mu \tilde{F}^{\mu\nu} = 0 \ . \tag{2.21}$$

By definition (2.14) the field tensor is invariant under a gauge transformation

$$A^\mu \longrightarrow A^{\mu\prime} = A^\mu + \partial^\mu \phi \ . \tag{2.22}$$

The same then holds for the equation of motion (2.19). This freedom can be used to impose constraints on the four components of the vector field A^μ.

2.1.1.2 Massive vector fields

Vector fields in which – in contrast to the electromagnetic field - the field quanta are massive are described by the so-called Proca equation

$$\partial_\mu F^{\mu\nu} + m^2 A^\nu = j^\nu \ . \tag{2.23}$$

Operating on this equation with the four-divergence ∂_ν gives, because F is antisymmetric,

$$m^2 \partial_\nu A^\nu = \partial_\nu j^\nu \ . \tag{2.24}$$

For $m \neq 0$ and a conserved current, this reduces the equation of motion (2.23) to

$$\left(\Box + m^2\right) A^\nu = j^\nu \ ; \qquad \partial_\nu A^\nu = 0 \ . \tag{2.25}$$

Thus for massive vector fields the freedom to make gauge transformations on the vector field is lost. The gauge-condition reduces the degrees of freedom of the field from 4 to 3. The space-like components represent the physical degrees of freedom.

The Lagrangian that leads to (2.23) is given by

$$\mathcal{L} = -\frac{1}{4} F^2 + \frac{1}{2} m^2 A^2 - j \cdot A \ . \tag{2.26}$$

2.1.1.3 Klein-Gordon fields

A particularly simple example is provided by the so-called Klein-Gordon field ϕ that obeys the equation of motion

$$\left(\partial_\mu \partial^\mu + m^2\right) \phi = \left(\Box + m^2\right) \phi = 0 \ ; \tag{2.27}$$

such a field describes scalar particles, i.e. particles without intrinsic spin. The Lagrangian leading to (2.27) is given by

$$\mathcal{L} = \frac{1}{2}\left((\partial_\mu \phi)(\partial^\mu \phi) - m^2\phi^2\right) \ . \tag{2.28}$$

An interesting case occurs if we consider 2 independent real scalar fields, ϕ_1 and ϕ_2, with the same mass m. The total Lagrangian is then simply given by a sum over the Lagrangians describing the individual fields, i.e.

$$\mathcal{L} = \frac{1}{2}\left((\partial_\mu \phi_1)(\partial^\mu \phi_1) - m^2\phi_1^2\right) + \frac{1}{2}\left((\partial_\mu \phi_2)(\partial^\mu \phi_2) - m^2\phi_2^2\right) \ . \tag{2.29}$$

On the other hand, we can also construct from the 2 real fields ϕ_1 and ϕ_2 2 complex fields, namely

$$\phi = \frac{1}{\sqrt{2}}(\phi_1 + \mathrm{i}\phi_2) \tag{2.30}$$

and its complex conjugate. In terms of these the Lagrangian (2.29) can be rewritten into

$$\mathcal{L} = (\partial_\mu \phi)^\dagger (\partial^\mu \phi) - m^2\phi^\dagger\phi \ . \tag{2.31}$$

2.1.1.4 Dirac Fields

A particularly simple example is provided by the Dirac field Φ for which the equation of motion is just the Dirac equation

$$(\mathrm{i}\gamma^\mu\partial_\mu - m)\,\Psi = 0 \ , \tag{2.32}$$

where the γ_μ are the usual (4×4) matrices of Dirac theory (see appendix A.1).

The corresponding Lagrangian is given by

$$\mathcal{L} = \bar{\Psi}\,(i\gamma^\mu\partial_\mu - m)\,\Psi \quad . \tag{2.33}$$

This can be seen by identifying the fields Φ_α in eq. (2.7) with the four components of the Dirac spinor $\bar{\Psi} = \Psi^\dagger\gamma_0$. Since $\mathcal{L}$ does not depend on $\partial_\mu\bar{\Psi}$ the equation of motion is simply given by

$$\frac{\partial\mathcal{L}}{\partial\bar{\Psi}} = (\mathrm{i}\gamma^\mu\partial_\mu - m)\,\Psi = 0 \tag{2.34}$$

2.2 Symmetries and Conservation Laws

As in classical mechanics there is also in field theory a conservation law associated with each continuous symmetry of $\mathcal{L}$. The theorem which describes the connection between the invariance of the Lagrangian under a continuous symmetry transformation and the related conserved current is known as Noether's theorem. In the following this will be illustrated for different types of symmetries which then lead to the well known conservation laws.

2.2.1 Geometrical Space-Time Symmetries

In this section we investigate the consequences of translations in four-dimensional space-time, i.e. infinitesimal transformations of the form

$$x_\nu \to x'_\nu = x_\nu + \epsilon_\nu \quad , \tag{2.35}$$

where the infinitesimal shift ϵ_ν is independent of the coordinate x_ν. Under such transformations the change of $\mathcal{L}$ is given by

$$\delta\mathcal{L} = \epsilon_\nu \frac{\partial \mathcal{L}}{dx_\nu} = \epsilon_\nu \partial^\nu \mathcal{L} \quad , \tag{2.36}$$

since $\mathcal{L}$ is a scalar.

If now $\mathcal{L}$ is required to be form-invariant under translations, it does not explicitly depend on x_ν, so that the derivative ∂_ν in (2.36) acts only on the arguments of $\Phi(x_\nu)$ and $\partial_\nu\Phi(x_\nu)$. The change of $\mathcal{L}$ can therefore also be expressed as

$$\delta\mathcal{L} = \frac{\partial \mathcal{L}}{\partial \Phi_\alpha}\delta\Phi_\alpha + \frac{\partial \mathcal{L}}{\partial(\partial_\mu \Phi_\alpha)}\delta(\partial_\mu \Phi_\alpha) \quad . \tag{2.37}$$

Since the variation and the differentiation are linear operations their order can be interchanged. We thus obtain

$$\delta\mathcal{L} = \frac{\partial \mathcal{L}}{\partial \Phi_\alpha}\delta\Phi_\alpha + \frac{\partial \mathcal{L}}{\partial(\partial_\mu \Phi_\alpha)}\partial_\mu\left(\delta \Phi_\alpha\right) \quad . \tag{2.38}$$

Using the equation of motion (2.7) for rewriting the first term yields

$$\begin{aligned} \delta\mathcal{L} &= \partial_\mu\left(\frac{\partial \mathcal{L}}{\partial(\partial_\mu \Phi_\alpha)}\right)\delta\Phi_\alpha + \frac{\partial \mathcal{L}}{\partial(\partial_\mu \Phi_\alpha)}\partial_\mu\left(\delta \Phi_\alpha\right) \\ &= \partial_\mu\left(\frac{\partial \mathcal{L}}{\partial(\partial_\mu \Phi_\alpha)}\delta\Phi_\alpha\right) \quad . \end{aligned} \tag{2.39}$$

The changes of the fields Φ_α appearing here are given by

$$\delta\Phi_\alpha(x) = \epsilon_\nu \frac{\partial \Phi_\alpha}{\partial x_\nu} = \epsilon_\nu \partial^\nu \Phi_\alpha(x) \tag{2.40}$$

Inserting eqs. (2.40) into (2.39) yields

$$\delta\mathcal{L} = \epsilon_\nu \partial_\mu \left(\frac{\partial \mathcal{L}}{\partial(\partial_\mu \Phi_\alpha)} \partial^\nu \Phi_\alpha \right) \quad . \tag{2.41}$$

Equating eqs. (2.41) and (2.36) finally gives

$$\partial_\mu \left(\frac{\partial \mathcal{L}}{\partial(\partial_\mu \Phi_\alpha)} \partial^\nu \Phi_\alpha - \mathcal{L}\, g^{\nu\mu} \right) = 0 \quad , \tag{2.42}$$

since the ϵ_ν are arbitrary. By defining the tensor $T^{\mu\nu}$ as

$$T^{\mu\nu} \equiv \frac{\partial \mathcal{L}}{\partial(\partial_\mu \Phi_\alpha)} \partial^\nu \Phi_\alpha - \mathcal{L}\, g^{\mu\nu} \tag{2.43}$$

this equation reads

$$\partial_\mu T^{\mu\nu} = 0 \quad . \tag{2.44}$$

Relation (2.44) has the form of a continuity equation. Spatial integration over a finite volume yields

$$\frac{d}{dt} \left(\int\limits_V T^{0\mu}(x) d^3x \right) = - \int\limits_V \frac{\partial T^{i\mu}}{\partial x_i}\, d^3x = \oint_S \vec{S}^{(\mu)} \cdot \vec{n}\, dS \quad . \tag{2.45}$$

Here $\vec{n}$ is a unit vector vertical on the surface S pointing outwards and $\vec{S}^{(\mu)}$ is a three-vector:

$$\vec{S}^{(\mu)} = -(T_{1\mu}, T_{2\mu}, T_{3\mu}) \quad . \tag{2.46}$$

The surface integral on the rhs of eq. (2.45) is taken over the surface S of volume V. For localized fields it can be made to vanish by extending the volume towards infinity. It is then evident that the quantities

$$P^\mu = \int T^{0\mu} d^3x \tag{2.47}$$

are conserved. These are the components of the four-momentum of the field, as can be verified for the zeroth component,

$$\begin{aligned} P^0 &= \int T^{00}\, d^3x = \int \left(\frac{\partial \mathcal{L}}{\partial(\partial_0 \Phi_\alpha)} \partial^0 \Phi_\alpha - \mathcal{L} \right) d^3x \\ &= \int (\Pi_\alpha \dot{\Phi}_\alpha - \mathcal{L})\, d^3x = H \quad , \end{aligned} \tag{2.48}$$

according to eqs. (2.9) and (2.10). The spatial components of the field momentum are

$$P^k = \int T^{0k}\, d^3x = \int \frac{\partial\mathcal{L}}{\partial(\partial_0\Phi_\alpha)}\partial^k\Phi_\alpha\, d^3x \quad . \tag{2.49}$$

In classical mechanics the form invariance of the Lagrangian under rotations leads to the conservation of angular momentum. Analogously, in a relativistic field theory the form invariance of $\mathcal{L}$ under four-dimensional space-time rotations (Lorentz covariance) leads to the conservation of a quantity that is identified with the angular momentum of the field. To obtain the same form for the angular momentum as in classical mechanics it is essential that $T_{\nu\mu}$ is symmetric. This can always be achieved, because Eq. (2.44) does not define the tensor T uniquely. We can always add a term of the form $\partial_\lambda D^{\lambda\mu\nu}$, where $D^{\lambda\mu\nu}$ is a tensor antisymmetric in the indices λ and μ, such that T becomes symmetric.

Comparing (2.45) with (2.49) and assuming T to be symmetric we see that the normal components of the vectors $\vec{S}^{(\mu)}$ in eq. (2.46) describe the momentum flow through the surface S of the volume V and thus determine the "radiation pressure" of the field[2]. These properties allow us to identify $T_{\mu\nu}$ as the energy-momentum tensor of the field. For the Lagrangian (2.17) of electrodynamics $T_{\mu\nu}$ is just the well-known Maxwell's stress tensor.

As already mentioned at the beginning of this chapter these conservation laws are special cases of Noether's theorem, which can be summarized for the general case as follows:

> *Each continuous symmetry transformation that leaves the Lagrangian invariant is associated with a conserved current. The spatial integral over this current's zeroth component yields a conserved charge.*

2.2.2 Internal Symmetries

Relativistic field theories may contain conservation laws that are not consequences of space-time symmetries of the Lagrangian, but instead are connected with symmetries in the internal degrees of freedom such as, e. g. , isospin or charge.

We therefore now allow for a mixture of the different fields under the transformation

$$\Phi_\alpha(x) \to \Phi'_\alpha(x) = \mathrm{e}^{-\mathrm{i}\epsilon q_{\alpha\beta}}\Phi_\beta \quad , \tag{2.50}$$

[2]More precisely, $S_k^{(\mu)}$ denotes the flux of the μ th component of the field momentum in the direction x_k.

where ϵ is an infinitesimal parameter and the $q_{\alpha\beta}$ are fixed c-numbers. We then have

$$\delta\Phi_\alpha(x) = \Phi'_\alpha(x) - \Phi_\alpha(x) = -\mathrm{i}\ \epsilon\ q_{\alpha\beta}\Phi_\beta(x) \quad . \tag{2.51}$$

The change of the Lagrangian is in general given by eq. (2.39)

$$\delta\mathcal{L} = \partial_\mu \left(\frac{\partial\mathcal{L}}{\partial(\partial_\mu\Phi_\alpha)} \delta\Phi_\alpha \right) \quad . \tag{2.52}$$

If $\mathcal{L}$ is invariant under this variation $\delta\Phi_\alpha$, then we have

$$\delta\mathcal{L} = \partial_\mu \left(\frac{\partial\mathcal{L}}{\partial(\partial_\mu\Phi_\alpha)} \delta\Phi_\alpha \right) = 0 \quad . \tag{2.53}$$

Eq. (2.53) is in the form of a continuity equation for the "current"

$$j^\mu(x) = \frac{\partial\mathcal{L}}{\partial(\partial_\mu\Phi_\alpha)} \delta\Phi_\alpha \frac{1}{\epsilon} \quad . \tag{2.54}$$

Inserting the field variations $\delta\Phi_\alpha$ yields for the current

$$j^\mu(x) = -\mathrm{i} \frac{\partial\mathcal{L}}{\partial(\partial_\mu\Phi_\alpha)} q_{\alpha\beta}\Phi_\beta \quad . \tag{2.55}$$

Eqs. (2.54) and (2.55) imply that the "charge"

$$Q = \int j^0(x)\, d^3x = -\mathrm{i} \int \frac{\partial\mathcal{L}}{\partial(\partial_0\Phi_\alpha)} q_{\alpha\beta}\Phi_\beta\, d^3x \tag{2.56}$$

of the system is conserved. The physical nature of these "charges" and "currents" has to remain open. It depends on the specific form of the symmetry transformation (2.50) and can be determined only by coupling the system to external fields.

Example: QED. To illustrate this conservation law, the theory of electromagnetic interactions is used as an example. However, in contrast to the considerations in section 2.1.1 we now consider a coupled system of a fermion field $\Psi(x)$ and the electromagnetic field $A_\mu(x)$ to determine the physical meaning of the conserved current. Together with a quantization procedure this theory is called Quantum Electrodynamics (QED). The Lagrangian is given by

$$\mathcal{L} = -\frac{1}{4} F_{\mu\nu}F^{\mu\nu} + \bar{\Psi}\left[\mathrm{i}\gamma^\mu(\partial_\mu + \mathrm{i}eA_\mu) - m\right]\Psi \quad . \tag{2.57}$$

$\mathcal{L}$ contains a part that describes the free electromagnetic field (first term). The second term, to be discussed and explained in detail later (in chapt. 7), describes the fermion Lagrangian; it is obtained from the free particle Lagrangian of eq. (2.33) by replacing the derivative ∂_μ by the covariant derivative

$$D_\mu = \partial_\mu + \mathrm{i}eA_\mu \tag{2.58}$$

(minimal coupling).

The Lagrangian (2.57) is obviously invariant under a variation of the fermion fields of the form

$$\Psi \to \Psi' = \mathrm{e}^{-\mathrm{i}\epsilon e}\Psi \quad . \tag{2.59}$$

Comparison with eq. (2.50) gives $q_{\alpha\beta} = e\ \delta_{\alpha\beta}$ so that the conserved "current" given by eq. (2.55) is:

$$j_\mu(x) = e\ \bar{\Psi}\gamma_\mu\Psi \quad . \tag{2.60}$$

Note that this conserved current is exactly the quantity that couples to the electromagnetic field in eq. (2.57). This property allows one to identify the current (2.60) as the electromagnetic current of the electron fields.

Example: Complex scalar field. The Lagrangian for the case of a complex scalar field interacting with an electromagnetic field is given by

$$\mathcal{L} = -\frac{1}{4}F_{\mu\nu}F^{\mu\nu} + (D_\mu\phi)^\dagger\,(D^\mu\phi) - m^2\phi^\dagger\phi\ . \tag{2.61}$$

This Lagrangian is simply the sum of the free electromagnetic Lagrangian (2.17) and the Lagrangian for a complex scalar field (2.31), where in the latter again the derivative ∂_μ has been replaced – through minimal substitution – by the covariant derivative D_μ (2.58).

The Lagrangian (2.61) is obviously invariant under the phase transformations

$$\begin{aligned} \phi(x) &\longrightarrow e^{-\mathrm{i}\varepsilon e}\phi(x) \\ \phi^\dagger(x) &\longrightarrow e^{+\mathrm{i}\varepsilon e}\phi^\dagger(x) \end{aligned} \tag{2.62}$$

The conserved current connected with this invariance can be obtained from the definition (2.55)

$$\begin{aligned} j^\mu &= -\mathrm{i}\left(\frac{\partial\mathcal{L}}{\partial\left(\partial_\mu\phi\right)}e\phi + \frac{\partial\mathcal{L}}{\partial\left(\partial_\mu\phi^\dagger\right)}(-e)\phi^\dagger\right) \\ &= \mathrm{i}e\left(\phi^\dagger D^\mu\phi - \phi\left(D^\mu\phi\right)^\dagger\right)\ . \end{aligned} \tag{2.63}$$

The conserved charge is then given by

$$Q = \int d^3x \, j^0(x) = \mathrm{i}e \int d^3x \, \left(\phi^\dagger D^0 \phi - \phi \left(D^0 \phi \right)^\dagger \right) \ . \tag{2.64}$$

It is evident that the charge disappears if the field ϕ is real. It is furthermore remarkable that now the electromagnetic field A^μ appears in the conserved current (through the covariant derivative $D^\mu = \partial^\mu + \mathrm{i}eA^\mu$). Again the conserved current provides the coupling to the electromagnetic field.

Part III

Global Symmetries

Chapter 3

SYMMETRIES OF MESON AND BARYON SYSTEMS

In chapter 2 it was shown how geometrical and internal symmetries of the Lagrangian lead to conserved currents and charges. In this chapter we take the opposite approach of using the observations of conservation laws to deduce the required symmetries of the Lagrangian. These, together with a desire for simplicity, will allow us to formulate a Lagrangian.

To demonstrate this construction of a Lagrangian out of the underlying symmetry properties, we first consider the physics of strongly interacting particles, namely mesons and baryons. In this context we consider both as elementary particles. This is a very useful approximation for describing nuclear phenomena at low and medium energies where any possible substructure could not be resolved, but is certainly not valid for the description of high-energy processes.

Nevertheless, historically this approach has lead to the prediction of the substructure of the mesons and baryons. This will be discussed later in this book. After some general remarks about the underlying symmetries, we will start out with the pion-nucleon system as the simplest example for the variety of the known systems of mesons and baryons.

3.1 $U(1)$ Symmetry

The simplest transformations of the fields appearing in a Lagrangian are global one-parameter changes of their phases. These transformations (2.50) with $q_{\alpha\beta} = q\delta_{\alpha\beta}$ for different values of the only free parameter ϵ form a group, called $U(1)$. Invariance under $U(1)$ transformations is connected with charge conservation (cf. the discussion at the end of chapter 2) (remember that the term "charge" does not necessarily denote the electric charge; the nature of the charge can be determined only by coupling it to an external field).

The physical conservation laws connected with invariance of the Lagrangian under these transformations are those for the electric charge and the baryon and lepton numbers.

The conservation of the total electric charge of a closed system can be concluded from the electric neutrality of macroscopic matter. This yields very stringent tests for the conservation of charge with time. Furthermore, in all known reactions charge is also conserved; for example, in the β-decay of the neutron one finds for the relative change of charge: $\Delta q/q \leq 10^{-23}$.

Particles and resonance states that decay ultimately into protons are called baryons. These particles all have half-integer spins and masses higher than that of the nucleon. The baryons can only decay into other lower-mass baryons, and cannot disappear altogether; at present the lower limit for the half-life of the ultimate decay product, the proton, is $T_{1/2} \geq 10^{32}$ years. This empirical conservation can be expressed by introducing a baryon number $B = 1$ for the baryons, $B = -1$ for the antibaryons and $B = 0$ for all other particles. B is then conserved in all processes known up today.

To finish our brief discussion of $U(1)$-symmetries observed in nature, we discuss the application to the six leptons known at present (e, μ, τ and their neutrinos) together with their antiparticles. Like the baryons, these have all half-integer spin but interact only through the weak and the electromagnetic forces. Again, one observes that, if one ascribes a lepton number $L = 1$ to each generation of leptons, i.e. to each lepton and its neutrino separately, and $L = -1$ to their antiparticles, then the lepton number L is a conserved quantity for each generation.

All of these conservation laws are consequences of $U(1)$ symmetries of the Lagrangian that will be discussed in this section.

3.1.1 Properties of the Group $U(1)$

In chapter 2 it was shown that there exists a conserved Noether current connected with an invariance of $\mathcal{L}$ under the global phase transformation[1]

$$\Phi(x) \rightarrow \Phi'(x) = \mathrm{e}^{-\mathrm{i}\epsilon q}\Phi(x) \quad , \tag{3.1}$$

[1]The following considerations hold for all types of fields, scalar, vector and Dirac.

where ϵ is an arbitrary constant and q a strength parameter for the conserved charge. The unitary transformations

$$U(\epsilon) = \mathrm{e}^{-\mathrm{i}\epsilon q} \quad , \tag{3.2}$$

with the fixed real constant q and labeled by the real parameter ϵ, form the elements of the so-called $U(1)$ group, the group of unitary transformations in one dimension (in general $U(n)$ is defined as the group of unitary matrices of dimension $n \times n$).

Before we start with the explicit construction of a $U(1)$-symmetric Lagrangian in the next section, we can already make some statements about its structure, based on the requirement of $U(1)$ invariance alone. A free Lagrangian that is required to lead to charge conservation can contain the fields only in such a form that it is invariant under the transformation (3.1). This can obviously be achieved only if $\mathcal{L}$ depends only on combinations of the form:

$$\bar{\Psi}(x) O \Psi(x) \tag{3.3a}$$

for fermions or

$$\Phi^\dagger O \Phi \tag{3.3b}$$

for bosons, where O is an operator acting in space and time or some internal space like, e.g. for the case of fermions, in the internal spinor space.

Quantummechanical consistency. It is instructive to consider whether after quantization the Noether charge $\hat{Q}$ is still a conserved quantity. Necessary for the quantum mechanical conservation law is that $\hat{Q}$ commutes with the Hamilton operator $\hat{H}$

$$\left[\hat{H}, \hat{Q}\right] = 0 \tag{3.4}$$

as follows directly from Heisenberg's equation of motion. Eq. (3.4) implies that the Hamiltonian operator $\hat{H}$ is invariant under the unitary transformation

$$\hat{H} \longrightarrow \hat{H}' = \mathrm{e}^{\mathrm{i}\epsilon\hat{Q}} \hat{H} \mathrm{e}^{-\mathrm{i}\epsilon\hat{Q}} \quad . \tag{3.5}$$

Since the classical Hamiltonian H of a physical system is always a $U(1)$-invariant functional of Φ and $\Phi^\dagger$ (see, e. g. eqs. (2.57) or (2.61)), the invariance (3.5) follows if the *field operators* transform as

$$\Phi \to \Phi' = \mathrm{e}^{\mathrm{i}\epsilon\hat{Q}} \Phi \mathrm{e}^{-\mathrm{i}\epsilon\hat{Q}} = \mathrm{e}^{-\mathrm{i}\epsilon q} \Phi \quad , \tag{3.6}$$

i. e. if the unitary transformation of Φ changes just the phase of Φ. Condition (3.6) leads immediately to

$$\left[\hat{Q}, \Phi\right] = -q\Phi \quad . \tag{3.7}$$

This relation shows that the consistency condition (3.6) implies that q has to be an eigenvalue of $\hat{Q}$ as can be seen by applying eq. (3.7) to the vacuum state

$$\begin{aligned} \hat{Q}\Phi|0> &= \Phi\hat{Q}|0> -q\Phi|0> \\ &= 0 - q\Phi|0> \quad . \end{aligned} \tag{3.8}$$

Essential for this result is the assumption that $\hat{Q}|0>= 0$, i.e. that the physical vacuum is uncharged.

Field quantization tells us that the state $\Phi|0>$ is indeed a one-antiparticle state; eq. (3.8) then tells us that the Noether charge is also conserved in a quantized theory if the parameter q appearing in the classical charge is an eigenvalue of the charge operator $\hat{Q}$. Turning this argument around, we could also postulate that the conservation law is quantummechanically consistent; this would then lead to eq. (3.8) and thus to the general result that every field operator creates an antiparticle and annihilates a particle. As an example we consider the charge operator for a charged scalar field ϕ

$$Q = \mathrm{i}q \int d^3x \, \left(\phi^\dagger \partial_0 \phi - \phi \partial_0 \phi^\dagger\right) \; . \tag{3.9}$$

By using equal-time commutation relations for the field operators we indeed find that (3.7) is fulfilled.

3.1.2 Structure of the Nucleon Lagrangian

As a first example we now construct a Lagrangian that describes free neutrons and protons as elementary particles. Since baryon-number is experimentally known to be conserved, we require the corresponding $U(1)_B$ symmetry from the outset. The Lagrangian is then that of free fermions for both the protons and the neutrons

$$\mathcal{L}_{pn} = \bar{\Psi}_p(\mathrm{i}\gamma^\mu\partial_\mu - M_p)\Psi_p + \bar{\Psi}_n(\mathrm{i}\gamma^\mu\partial_\mu - M_n)\Psi_n \quad . \tag{3.10}$$

The experimental values for the masses of proton and neutron are $M_p = 938.3$ MeV and $M_n = 939.6$ MeV, respectively.

$\mathcal{L}_{pn}$ is obviously invariant under the $U(1)$ transformation

$$\begin{array}{lll} \Psi_p & \to & \Psi_p' = \mathrm{e}^{-\mathrm{i}\epsilon q_p}\Psi_p, \qquad \bar{\Psi}_p \to \bar{\Psi}_p' = \mathrm{e}^{+\mathrm{i}\epsilon q_p}\bar{\Psi}_p \\ \Psi_n & \to & \Psi_n' = \mathrm{e}^{-\mathrm{i}\epsilon q_n}\Psi_n, \qquad \bar{\Psi}_n \to \bar{\Psi}_n' = \mathrm{e}^{+\mathrm{i}\epsilon q_p}\bar{\Psi}_n \quad , \end{array} \tag{3.11}$$

since it depends only on bilinear field combinations of the form $\bar{\Psi}\Psi$. Assigning the value $q_n = q_p = 1$ to both the proton and neutron charge leads to conservation of the baryon number

$$B = \int \left(\Psi_p^\dagger \Psi_p + \Psi_n^\dagger \Psi_n\right) d^3x \quad ; \tag{3.12}$$

the corresponding symmetry operation is $U(1)_B$.

In addition to baryon number conservation, charge conservation also holds; in this case one chooses $q_p = 1$, $q_n = 0$, so that the transformation is applied to the proton field alone. The corresponding transformations

$$\begin{array}{lll} \Psi_p & \to & \Psi_p' = \mathrm{e}^{-\mathrm{i}\epsilon}\Psi_p, \qquad \bar{\Psi}_p \to \bar{\Psi}_p' = \mathrm{e}^{+\mathrm{i}\epsilon}\bar{\Psi}_p \\ \Psi_n & \to & \Psi_n' = \Psi_n, \qquad \bar{\Psi}_n \to \bar{\Psi}_n' = \bar{\Psi}_n \end{array}$$

form another group, $U(1)_Q$, with the conserved charge

$$Q = \int \Psi_p^\dagger \Psi_p \, d^3x \quad . \tag{3.13}$$

Both of these conservation laws hold in nature. Another one, that could be obtained by setting $q_p = 0$, $q_n = 1$, would lead to neutron number conservation that, however, is violated in nature.

The close agreement of the experimental values of neutron and proton masses suggests to view the small difference between them as the consequence of a perturbative interaction. Abbreviating the latter as $\mathcal{L}_{\text{int}}$ allows us to rewrite the Lagrangian (3.10) as

$$\mathcal{L}_{pn} = \sum_{q=p,n} \left[\bar{\Psi}_q \left(\mathrm{i}\gamma^\mu \partial_\mu - M\right) \Psi_q\right] + \mathcal{L}_{\text{int}} \quad , \tag{3.14}$$

where M is the common mass for proton and neutron in the absence of the perturbing interaction.

We now formally introduce a two-dimensional spinor

$$\Psi_N = \begin{pmatrix} \Psi_p \\ \Psi_n \end{pmatrix} \quad , \tag{3.15}$$

in terms of which the Lagrangian (3.14) can be written as

$$\begin{aligned}\mathcal{L}_{pn} &= \bar{\Psi}_N \left(\mathrm{i}\gamma^\mu \partial_\mu - M\right) \Psi_N + \mathcal{L}_{int} \\ &= \mathcal{L}_N + \mathcal{L}_{int} \quad .\end{aligned} \tag{3.16}$$

Here the Lagrangian has been split into two parts; $\mathcal{L}_N$ is that of the noninteracting, equal-mass protons and neutrons whereas $\mathcal{L}_{int}$ denotes the residual interaction responsible for the small mass difference between proton and neutron.

The form (3.15) for the state Ψ_N suggests to view proton and neutron as two different intrinsic states of the same particle, the nucleon, described by the Lagrangian $\mathcal{L}_N$. The new degree of freedom that distinguishes between proton and neutron is called isospin.

3.2 $SU(2)$ Symmetry

From eq. (3.14) it is evident that $\mathcal{L}_{pn}$ is invariant under an exchange of protons and neutrons, if the relatively small $\mathcal{L}_{int}$ is neglected. That the total strong interaction Lagrangian is indeed invariant under such a transformation is suggested by experimental observations, for example, the near-equality of the binding energy and spectra of the mirror nuclei ^{7}Li and ^{7}Be that differ by the change of one neutron into one proton. The structural similarity of these two nuclei indicates that proton-proton and neutron-neutron forces are nearly equal. Thus the strong interaction Lagrangian seems to be invariant under an exchange of protons and neutrons. Therefore, it can also not distinguish between any mixture of protons and neutrons. This implies that $\mathcal{L}_N$ of eq. (3.16) contains a further symmetry – beyond the simple $U(1)$ – that mixes the field components Ψ_p and Ψ_n. To discuss this symmetry is the purpose of this section.

3.2.1 Properties of the Group $SU(2)$

We consider a space- and time-independent general transformation of the two-dimensional spinor Ψ_N which was defined in (3.15)[2]

$$\Psi \to \Psi' = U\Psi \quad . \tag{3.17}$$

In its most general form U must be a (2×2)-dimensional matrix with space- and time-independent complex components. We will now look for such a transformation U that leaves $\mathcal{L}_N$ invariant.

[2]We suppress here, as in the following, the subscript N of Ψ.

The fields Ψ' must also fulfill the anticommutation relations for fermions. This requirement implies that the transformation U must be unitary:

$$U^\dagger U = UU^\dagger = 1 \quad . \tag{3.18}$$

Because of this property one obtains for the determinant of the transformation matrix U

$$\det(U^\dagger)\det(U) = (\det(U))^* \det(U) = 1 \tag{3.19}$$

and thus

$$\det(U) = \mathrm{e}^{-\mathrm{i}\varphi} \qquad (\varphi \text{ real}) \quad . \tag{3.20}$$

Eq. (3.20) implies that U contains a simple $U(1)$ transformation. Since the $U(1)$ symmetry has also already been incorporated we can disregard it here by simply setting $\varphi = 0$. Then one has

$$\det(U) = 1 \quad . \tag{3.21}$$

Conditions (3.18) and (3.20) narrow the possible matrices U down to the unitary matrices with unit determinant, the so-called special unitary matrices. The set of all of these (2×2)-matrices forms a group, the Lie group $SU(2)$; the product of any two of the elements of this group is just the normal matrix multiplication.

Lie groups have the special property that their elements can be parametrized smoothly and analytically such that each element can be obtained by a continuous parameter change from the identity. In order to obtain the parametrization of the group elements it is, therefore, sufficient to consider an infinitesimal change of the spinor Ψ

$$\Psi' = (1 - i\xi)\Psi \tag{3.22}$$

where ξ is a (2×2)-dimensional matrix with infinitesimally small matrix elements and 1 is the (2×2)-dimensional unit matrix. The unitarity condition (3.18) then reads

$$\begin{aligned} U^\dagger U &= (1 + \mathrm{i}\xi^\dagger)(1 - \mathrm{i}\xi) \\ &= 1 + \mathrm{i}(\xi^\dagger - \xi) + \mathcal{O}(\xi^2) = 1 \quad . \end{aligned} \tag{3.23}$$

This can be fulfilled only if

$$\xi^\dagger = \xi \quad , \tag{3.24}$$

i.e. if ξ is a hermitian matrix. Furthermore, from

$$1 = \det(U) = 1 - \mathrm{i}\xi_{11} - \mathrm{i}\xi_{22} + \mathcal{O}(\xi^2) \tag{3.25}$$

follows

$$\mathrm{tr}(\xi) = 0 \quad . \tag{3.26}$$

Equations (3.24) and (3.26) show that ξ is a hermitean matrix with vanishing trace. Since its dimension is (2×2), it contains three independent elements. Such matrices can in general be written as a linear superposition of three linearly independent, hermitian, traceless basis-matrices τ_i $(i = 1, 2, 3)$

$$\xi = \boldsymbol{\epsilon} \cdot \frac{\boldsymbol{\tau}}{2} \tag{3.27}$$

with the real parameters

$$\boldsymbol{\epsilon} = (\epsilon_1, \epsilon_2, \epsilon_3)$$

and the matrices

$$\boldsymbol{\tau} = (\tau_1, \tau_2, \tau_3) \quad .$$

Here the ϵ_i are infinitesimal, real constants and the τ_i are the Pauli matrices

$$\tau_1 = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \ , \quad \tau_2 = \begin{pmatrix} 0 & -\mathrm{i} \\ \mathrm{i} & 0 \end{pmatrix} \ , \quad \tau_3 = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} \quad . \tag{3.28}$$

These obey the commutation relations

$$\left[\frac{\tau_i}{2}, \frac{\tau_j}{2}\right] = \mathrm{i}\epsilon_{ijk}\frac{\tau_k}{2} \tag{3.29}$$

where ϵ_{ijk} is the Levi-Civita tensor:

$$\epsilon_{ijk} = \begin{cases} 1 & i,j,k = 1,2,3 \quad \text{and cyclical} \\ -1 & i,j,k = 3,2,1 \quad \text{and cyclical} \\ 0 & \text{in all other cases} \end{cases} \quad . \tag{3.30}$$

The commutation relations (3.29) define the "algebra" of the group $SU(2)$, the elements of the tensor ϵ_{ijk} are called the "structure constants" of $SU(2)$ and the matrices $\tau_i/2$ are the matrix representations of the "generators" of $SU(2)$-transformations in isospin space, i.e. in the space spanned by the vectors Ψ_N (eq. (3.15)). They are normalized such that

$$\mathrm{tr}(\tau_l \tau_k) = 2\delta_{lk} \ . \tag{3.31}$$

Summarizing all these results, the most general infinitesimal symmetry transformation of the two component spinor Ψ is given by

$$\Psi \to \Psi' = \left(1 - \mathrm{i}\boldsymbol{\epsilon} \cdot \frac{\boldsymbol{\tau}}{2}\right) \Psi \quad .$$

For a finite parameter $\boldsymbol{\epsilon}$ this is easily generalized to

$$\Psi \to \Psi' = \mathrm{e}^{-\mathrm{i}\boldsymbol{\epsilon}\frac{\boldsymbol{\tau}}{2}}\Psi \quad . \tag{3.32}$$

Formally, this is a generalization of the global gauge transformation of the $U(1)$ type with, however, two major differences:

1. There are now 3 "phase angles" ϵ_i .

2. The exponent in (3.32) now contains three matrices τ_i that do not commute.

Thus, two $SU(2)$ transformations will in general not commute. Consequently $SU(2)$ is a non-Abelian group. The special ($\det = 1$) unitary (2×2) matrices

$$U = \mathrm{e}^{-\mathrm{i}\boldsymbol{\epsilon t}} \quad , \qquad \text{with} \quad \boldsymbol{t} = \boldsymbol{\tau}/2 \tag{3.33}$$

form the fundamental representation of the group $SU(2)$; the matrices t_i are the matrix representations of the generators. The matrix U acts on two-dimensional vectors, the isospinors.

Higher-dimensional representations. There exist also matrices of higher dimension than (2×2) that obey the same $SU(2)$ algebra (3.29). These form higher dimensional representations of the symmetry group. For example, in 3 dimensions such matrices are

$$\begin{aligned} t_1 &= \frac{1}{\sqrt{2}}\begin{pmatrix} 0 & 1 & 0 \\ 1 & 0 & 1 \\ 0 & 1 & 0 \end{pmatrix} \quad t_2 = \frac{1}{\sqrt{2}}\begin{pmatrix} 0 & -\mathrm{i} & 0 \\ \mathrm{i} & 0 & -\mathrm{i} \\ 0 & \mathrm{i} & 0 \end{pmatrix} \\ t_3 &= \begin{pmatrix} 1 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & -1 \end{pmatrix} \end{aligned} \tag{3.34}$$

in a representation in which t_3 is diagonal. The elements of $SU(2)$ are then again given by eq. (3.33), with the t_i now defined by eq. (3.34).

Note that the generators t_i are not unique. All matrices that fulfill the commutation relations (3.29) are generators of $SU(2)$, though in a different representation. For example, the (3×3) matrices

$$(t_k)_{lm} = -\mathrm{i}\epsilon_{klm} \tag{3.35}$$

also obey the same commutation rules[3]. In the case of (3.35) the structure constants of $SU(2)$ define the matrix representation of the generators.

The $SU(2)$ transformation in a three-dimensional representation also acts on three-dimensional basis vectors. If we denote a vector in this space by $\mathbf{\Pi}$, then an infinitesimal transformation in the representation (3.35) is given by

$$\delta \Pi_l = -\mathrm{i}\,(\epsilon_k t_k \mathbf{\Pi})_l = -\epsilon_k \epsilon_{klm} \Pi_m = (\boldsymbol{\epsilon} \times \mathbf{\Pi})_l \quad , \tag{3.36}$$

i.e. just by the transformation law of an ordinary vector under an infinitesimal spatial rotation. This is a consequence of the fact that the algebra (3.29) for the $SU(2)$ generators $\tau_i/2$ is exactly the same as that for the orbital angular momentum operators l_i. The latter are the generators of the rotation group in three dimensions, $SO(3)$, that act on three-dimensional ordinary vectors. The two groups are thus locally equivalent.

General definitions. In general, the algebra of a group is defined by the commutation relation of its generators

$$[Q_i, Q_j] = \mathrm{i} c_{ijk} Q_k$$

where the c_{ijk} are the structure constants of the group. Then the matrices q_k with

$$(q_k)_{lm} = -\mathrm{i} c_{klm}$$

are the generators of the "regular" or "adjoint" representation of the group. Thus (3.35) defines the regular representation of $SU(2)$. The dimension of this special representation is equal to the number of group generators, i.e. $d = 3$ for $SU(2)$.

The representations (3.34) and (3.35) are examples of so-called "irreducible" representations that cannot be put into block-diagonal form by a similarity transformation, i.e. by a transformation $U \to U' = SUS^{-1}$ where S is any nonsingular, invertible matrix acting on the basis vectors of U. If this is possible for a given representation, then it is said to be reducible.

In general, in an n-dimensional representation, the group elements U act on an n-dimensional vector space with n linearly independent basis vectors. If the group representation is irreducible, then these basis vectors are said to form a "multiplet" . If the Hamiltonian is invariant under the transformation U, then all states of a multiplet have the same energy; they differ by the

[3]This can be shown by using the relation $\epsilon_{ijk}\epsilon_{klm} = \delta_{il}\delta_{jm} - \delta_{im}\delta_{jl}$.

eigenvalues of the diagonal generator, for example, by their value of t_3 in the irreducible representations (3.28) or (3.34).

Another important property of a Lie group is its rank, defined as the maximum number of simultaneously diagonal generators. In the set of the three Pauli matrices τ_i each matrix commutes only with itself; thus $SU(2)$ is a group of rank 1. From the members of an algebra of a Lie group of rank n one can, by combining the generators in a nonlinear form, construct n invariant operators, the so-called Casimir operators, which commute with all generators. For $SU(2)$, this Casimir operator is given by $\boldsymbol{\tau}^2 = \tau_1^2 + \tau_2^2 + \tau_3^2$. The Casimir operator is usually used to label the multiplets.

3.2.2 Application to the Pion-Nucleon System

The application of the considerations in the preceding section to physical particles is obtained if one can identify n particles with an n-dimensional multiplet of basis states. For example, in a theory with an $SU(2)$ invariant Hamiltonian this identification is possible if the n particles have approximately the same mass.

In this section we will now apply these symmetry considerations to the pion-nucleon system since this is one of the elementary and, at the same time, most relevant systems of nuclear physics. As already mentioned, the $SU(2)$ isospin symmetry is most evident in the charge independence of nuclear forces that can, for example, be observed in the binding energy of mirror nuclei. However, as will become clear from the discussions in this chapter, isospin symmetry goes further than charge independence, which is merely a consequence of it.

3.2.2.1 Symmetry of the Nucleon Lagrangian

We now apply some of the formal considerations of the last section to the case of the two-component nucleon spinor Ψ. Since

$$U = \mathrm{e}^{-\mathrm{i}\boldsymbol{\epsilon}\frac{\boldsymbol{\tau}}{2}} \tag{3.37}$$

is unitary, it is evident that the bilinear product $\bar{\Psi}\Psi$ in eq. (3.16) and thus the whole nucleon Lagrangian $\mathcal{L}_N$ is invariant under the $SU(2)$ transformation (3.37). The transformation U acts on the intrinsic degree of freedom that distinguishes between the two elements of the spinor Ψ, i.e. between proton and neutron. This internal degree of freedom is called isospin. The vectors $\boldsymbol{\epsilon}$ and $\boldsymbol{\tau}$ are then vectors in the isospin-space; they are called isovectors.

The conserved current corresponding to this $SU(2)$ symmetry can be obtained from expression (2.55) by generalizing it to the case, where $\boldsymbol{\epsilon}$ and $\boldsymbol{q}_{\alpha\beta}$ are vectors in the space of the internal degrees of freedom,

$$\boldsymbol{q}_{\alpha\beta} = \frac{\boldsymbol{\tau}_{\alpha\beta}}{2} \quad .$$

This yields for the conserved (isospin) current

$$\boldsymbol{j}_\mu = \bar{\Psi}\gamma_\mu \frac{\boldsymbol{\tau}}{2}\Psi \tag{3.38}$$

where $\boldsymbol{j}_\mu$ is now the μ th component of a four-vector in coordinate space and at the same time an isovector. The conserved charge is then another isovector,

$$\boldsymbol{T} = \int \boldsymbol{j}_0(x)\, d^3x = \int \Psi^\dagger(x)\frac{\boldsymbol{\tau}}{2}\Psi(x)\, d^3x \quad , \tag{3.39}$$

the *isospin*; Ψ is the two-dimensional spinor (3.15). Its third component can be written as

$$T_3 = \frac{1}{2}\int \left(\Psi_p^\dagger\Psi_p - \Psi_n^\dagger\Psi_n\right)\, d^3x \quad . \tag{3.40}$$

$\mathcal{L}_N$ is also invariant under the $U(1)_B$ transformation (3.11)

$$\Psi \to \Psi' = \mathrm{e}^{-\mathrm{i}\epsilon}\Psi \quad , \tag{3.41}$$

leading to the conservation of baryon number

$$B = \int \left(\Psi_p^\dagger\Psi_p + \Psi_n^\dagger\Psi_n\right)\, d^3x = \int \Psi^\dagger\Psi\, d^3x \quad . \tag{3.42}$$

Combining the symmetry operations (3.37) and (3.41), we see that there is also an invariance of $\mathcal{L}_N$ under the combined $U(1)_B \otimes SU(2)$ transformation

$$\Psi \to \Psi' = \mathrm{e}^{-\mathrm{i}\epsilon\frac{(1+\tau_3)}{2}}\Psi \quad , \tag{3.43}$$

leading to the conservation of the charge

$$Q = \int \Psi^\dagger \frac{1+\tau_3}{2}\Psi\, d^3x = \int \Psi_p^\dagger\Psi_p\, d^3x \quad . \tag{3.44}$$

Q is obviously equal to the number of electrical charges of the nucleons. Combining eqs (3.40), (3.42) and (3.44) yields the relation

$$Q = \frac{1}{2}B + T_3 \quad . \tag{3.45}$$

Although we have derived these relations only for the free nucleon Lagrangian (3.16), the empirically well founded isospin conservation together with that for electric charge and baryon number indicates that the full Lagrangian for the strong interaction respects these symmetries. Thus (3.45) must hold for any system of nucleons with $U(1)$ and $SU(2)$ symmetries if the Lagrangian contains no derivative terms other than those in the kinetic energy (in this case the currents remain unchanged); it is based only on the symmetry operations (3.37), (3.41) and (3.43) and not on any specific symmetry properties of the Lagrangian.

To determine the isospin T (defined via the eigenvalue $T(T+1)$ of the operator $\boldsymbol{T}^2$, which is just the Casimir operator of $SU(2)$) of a multiplet, we make use of the similarity of the group $SU(2)$ and the group of spatial rotations in three dimensions, $SO(3)$, whose generators are the angular momentum operators; both groups have the same algebra. A well known example for an $SO(3)$ symmetry is the hydrogen atom. There the spectrum consists of $2L+1$ dimensional multiplets labeled by the eigenvalues of the Casimir operator L^2; the states in each multiplet are labeled by L_3. Analogously, the n particles of the isospin multiplet have isospin T given via $n = 2T+1$ and can be distinguished via T_3.

In the case treated in this section proton and neutron are both members of the $T = 1/2$ multiplet, distinguished by $T_3 = +1/2$ and $T_3 = -1/2$, respectively. In an $SU(2)$-symmetric world they are thus two different states of the same particle, the nucleon.

The fundamental doublet of $SU(2)$ has $T = 1/2$ and $T_3 = \pm 1/2$. The corresponding antiparticles also form a doublet. Since the charge-conjugate spinor is given by

$$\Psi^C = \mathrm{i}\gamma^2\gamma^0\Psi^* \; , \tag{3.46}$$

the antiparticle spinor transforms under U^*. In $SU(2)$ we have

$$U^* = (-\mathrm{i}\tau_2)U(\mathrm{i}\tau_2) \tag{3.47}$$

because of $\vec{\tau}^* = -\tau_2\vec{\tau}\tau_2$. Note that the transformation U^* can be obtained from U by a similarity transformation; U and U^* are thus equivalent representations[4]. Therefore, we have

$$\Psi^{c\prime} = U^*\Psi^c = (-\mathrm{i}\tau_2)U(\mathrm{i}\tau_2)\Psi^c \tag{3.48}$$

and thus

$$(\mathrm{i}\tau_2\Psi^c)' = U(\mathrm{i}\tau_2\Psi^c) \tag{3.49}$$

[4] "Equivalent" matrix representations of $SU(N)$ can be transformed into each other by means of a unitary transformation belonging to the same group.

Thus, with

$$\Psi = \begin{pmatrix} \Psi_p \\ \Psi_n \end{pmatrix} \quad , \tag{3.50}$$

also

$$\Psi^C = (-\mathrm{i}\tau_2) \begin{pmatrix} \Psi_p^C \\ \Psi_n^C \end{pmatrix} = \begin{pmatrix} -\Psi_n^C \\ \Psi_p^C \end{pmatrix} \tag{3.51}$$

forms an $SU(2)$ doublet.

In analogy to the nucleon the particles that transform under the representation (3.34) have isospin $T = 1$ and $T_3 = +1, 0, -1$. If this representation is realized in nature then we should find a triplet of particles with the same mass and different electric charges. The $SU(2)$ transformation U, when applied to such a triplet, simply amounts to a rotation of a vector in isospin space.

In order to construct many-particle states with good $SU(2)$ transformation properties, the single-particle states of $SU(2)$ multiplets can also be coupled to states of different multiplets (i.e. different isospin) by using the normal Clebsch-Gordan coefficients well known from angular-momentum coupling. This analogy arises again because the group $SO(3)$ has the same algebra as $SU(2)$. Thus, e.g., the $T = 1$ multiplet mentioned at the start of this chapter could be obtained by coupling two nucleons each with $T = 1/2$.

Quantummechanical consistency. The components $\hat{T}_i$ of the isospin *operator*, after field quantization, obey the same commutation relations as the Pauli-matrices $\tau_i/2$

$$\left[\hat{T}_i, \hat{T}_j\right] = \mathrm{i}\epsilon_{ijk}\hat{T}_k \quad . \tag{3.52}$$

This follows from the general relation

$$\begin{aligned} &\left[\int \Psi^\dagger(x)\hat{O}\Psi(x)\, d^3x \;\; , \;\; \int \Psi^\dagger(y)\hat{Q}\Psi(y) d^3y \right] \\ &\qquad = \int \Psi^\dagger(x) \left[\hat{O}\, , \hat{Q}\right] \Psi(x)\, d^3x \qquad (x_0 = y_0) \end{aligned} \tag{3.53}$$

where all the points x and y are taken at the same time and $\hat{O}$ and $\hat{Q}$ are operators acting on the spinor components. Equation (3.53) can be easily proven if canonical *anti*commutation relations for the fields Ψ and $\Psi^\dagger$

are used. Considering the group properties, the operators $\hat{T}_i$ are therefore equivalent to the Pauli matrices

$$t_i = \frac{\tau_i}{2} \quad .$$

They only act onto different basis states, the $\hat{T}_i$ on Hilbert space state vectors and the t_i on basis vectors in the space of the intrinsic isospin degree of freedom.

Generalizing relation (3.7) to the isospin-case ($\hat{\boldsymbol{O}} = \boldsymbol{\tau}/2$) gives

$$\left[\hat{\boldsymbol{T}}, \Psi(x)\right] = -\frac{\boldsymbol{\tau}}{2}\Psi(x) \quad . \tag{3.54}$$

This is nothing other than the relation

$$\mathrm{e}^{\mathrm{i}\boldsymbol{\epsilon}\hat{\boldsymbol{T}}}\Psi\mathrm{e}^{-\mathrm{i}\boldsymbol{\epsilon}\hat{\boldsymbol{T}}} = \mathrm{e}^{-\mathrm{i}\boldsymbol{\epsilon}\frac{\boldsymbol{\tau}}{2}}\Psi \tag{3.55}$$

for infinitesimal $\boldsymbol{\epsilon}$. The lhs of this equation is the transformed field operator. Thus $\hat{\boldsymbol{T}}$, the conserved "charge" operator, is in analogy to the $U(1)$ case, the generator of the $SU(2)$ representation acting in Hilbert space. This again shows that the correspondence principle is satisfied, since the field operator transforms in the same way as the classical field. Also, as in the $U(1)$ case (eq. (3.8)), eq. (3.54) implies that $\tau_i/2$ is the eigenvalue of $\hat{T}_i$

$$\hat{T}_i\Psi|0> - \Psi\hat{T}_i|0> = -\frac{\tau_i}{2}\Psi|0> \quad . \tag{3.56}$$

Essential for this result is again the assumption that the vacuum is "uncharged", i.e. carries zero isospin. Formally, this corresponds to the use of a normal product in the definition of the charge in eq. (3.39), which here has not been written explicitly in order to simplify the notation.

Eq. (3.55) implies that the classical and the quantum symmetry operations have the same effect. Thus, if the Hamiltonian $\hat{H}$ of the system is invariant under (3.55), then the $\hat{T}_i$ represent conserved quantities also after quantization. This implies that $\hat{H}$ commutes with all the $\hat{T}_i$; its $SU(2)$-transformed eigenstates then belong to the same energy as the original ones.

3.2.2.2 Symmetry of the Pion Lagrangian

The pion carries the quantum numbers $J^\pi = 0^-$ and is thus a pseudoscalar particle. It exists in three different charge states, with nearly equal mass ($m(\pi^+) = m(\pi^-) = 139.6$ MeV, $m(\pi^0) = 135.0$ MeV); it is natural to associate the somewhat higher mass of the two charged pions with the effect of the repulsive Coulomb interaction. We therefore assume that in a world in which only the strong π–N force acts the three pions would have equal mass. This suggests that the three pions can be assigned to the three states of a $T = 1$ triplet, the basis vectors of the irreducible representation (3.34) and that, correspondingly, the strong interaction Lagrangian is symmetric under $SU(2)$ transformations. The pions thus have to be represented by an isovector with components π^+, π^0 and π^-; these are charged, nonhermitean pseudoscalar fields. That these components indeed belong to the representation (3.34) we can see by inspecting the eigenvalues of t_3 in that representation.

From these three components we can construct the basis triplet of a cartesian representation

$$\boldsymbol{\pi} = (\pi_1, \pi_2, \pi_3) \tag{3.57}$$

where the three components π_1, π_2 and π_3 are hermitean, pseudoscalar fields. These are obtained from the physical fields by means of the transformation

$$\begin{aligned}
\pi_1(x) &= \frac{1}{\sqrt{2}}(\pi^+(x) + \pi^-(x)) \\
\pi_2(x) &= \frac{\mathrm{i}}{\sqrt{2}}(\pi^+(x) - \pi^-(x)) \\
\pi_3(x) &= \pi^0(x) \quad .
\end{aligned} \tag{3.58}$$

The transformation (3.58) is exactly the same as the one that converts the components of an irreducible spherical tensor of rank 1, an object with well defined transformation properties under spatial rotations, into those of a normal, cartesian vector. This should come as no surprise since, as already mentioned at the end of the last subsection, $SO(3)$ has the same algebra as $SU(2)$.

The Lagrangian for the pions is given by that for charged scalar fields (cf. eq. (2.31))

$$\mathcal{L}_\pi = \left((\partial^\mu \pi^+)(\partial_\mu \pi^-) - m^2 \pi^+ \pi^-\right) + \frac{1}{2}\left((\partial_\mu \pi^0)^2 - m^2 (\pi^0)^2\right) \tag{3.59}$$

according to the general considerations in section 2.1.1.3. In terms of the cartesian isovector $\boldsymbol{\pi}$ (3.58) this can be rewritten as

$$\mathcal{L}_\pi = \frac{1}{2}\left[(\partial_\mu \boldsymbol{\pi})^2 - m^2 \boldsymbol{\pi}^2\right] \quad . \tag{3.60}$$

Written in this form, it is evident that $\mathcal{L}_\pi$ is $SU(2)$ symmetric since it contains only scalar products of the isovector $\boldsymbol{\pi}$ which are invariant under rotations in isospin space (eqs. (3.33), (3.37)).

The conserved current is

$$\boldsymbol{j}_\mu = -\mathrm{i}\,\partial_\mu \pi_k\,(\boldsymbol{t})_{kl}\,\pi_l = \boldsymbol{\pi} \times \partial_\mu \boldsymbol{\pi} \tag{3.61}$$

and the conserved charge is given by

$$\boldsymbol{T} = -\mathrm{i}\int \partial_t \pi_k\,(\boldsymbol{t})_{kl}\,\pi_l\,d^3x = \int (\boldsymbol{\pi} \times \boldsymbol{\Pi})\,d^3x \quad ; \tag{3.62}$$

$\boldsymbol{j}_\mu$ and $\boldsymbol{T}$ are in the cartesian, regular representation, both are vectors in isospin-space; $\boldsymbol{\Pi}$ is the conjugate field momentum.

The third component of $\boldsymbol{T}$ expressed by the multiplet states (3.57) is obtained as

$$T_3 = \mathrm{i}\int \left[(\partial_t \pi^+)\pi^- - (\partial_t \pi^-)\pi^+\right] d^3x \quad . \tag{3.63}$$

For the charge of the pions one has, from the $U(1)$ symmetry of (3.59),

$$Q = \mathrm{i}\int \left[(\partial_t \pi^+)\pi^- - (\partial_t \pi^-)\pi^+\right] d^3x \quad . \tag{3.64}$$

Comparing (3.64) and (3.63) gives

$$Q = T_3 \quad . \tag{3.65}$$

This relation can be viewed as a special case of eq. (3.45), since the pion has baryon number $B = 0$.

3.2.2.3 Symmetry of the Interaction

In this section we will now construct the Lagrangian for the pion-nucleon interaction, $\mathcal{L}_{\pi N}$. The first requirement of baryon number conservation means that $\mathcal{L}_{\pi N}$ can contain only bilinear forms in the nucleon field Ψ; in lowest order the coupling thus must read

$$\mathcal{L}_{\pi N} \sim \bar{\Psi}\hat{O}\Psi\,\hat{Q}\pi \quad , \tag{3.66}$$

where $\hat{O}$ is an operator that acts on the components of Ψ, and $\hat{Q}$ acts on the pion field. The latter can appear linearly in (3.66) since the pions carry no baryon number, so that single pions can be created or annihilated.

The structure of $\hat{O}$ can be obtained most easily by considering first interactions with the neutral pions only. Since π^0 is a pseudoscalar field (spin $= 0$, parity $= -1$) and the Lagrangian has to be a true scalar, $\bar{\Psi}\hat{O}\Psi$ itself must be a pseudoscalar, if $\hat{Q}$ is taken to be equal to 1 for simplicity. This can be achieved by an ansatz of the form (see eq. (A.31))[5]

$$\hat{O} = \gamma_5 \qquad \text{(pseudoscalar coupling)} \quad . \tag{3.67}$$

With this coupling, we obtain

$$\mathcal{L}_{\pi N} = \mathrm{i}g_1 \bar{\Psi}_n \gamma_5 \Psi_n \pi^0 + \mathrm{i}g_2 \bar{\Psi}_p \gamma_5 \Psi_p \pi^0 \tag{3.68}$$

where g_i is a real coupling strength. The factor i is necessary to make $\mathcal{L}_{\pi N}$ hermitian:

$$\left(\bar{\Psi}_n \gamma_5 \Psi_n \pi^0\right)^\dagger = \pi^0 \Psi_n^\dagger \gamma_5 \bar{\Psi}_n^\dagger = \pi^0 \Psi_n^\dagger \gamma_5 \gamma_0 \Psi_n = -\pi^0 \bar{\Psi}_n \gamma_5 \Psi_n \quad . \tag{3.69}$$

These considerations can be extended to the interaction with the charged pions. The requirement of both baryon number conservation and also charge conservation imposes immediately the constraint on $\mathcal{L}_{\pi N}$ that it has to contain combinations of the form

$$\bar{\Psi}_n \gamma_5 \Psi_p \pi^- \quad \text{or} \quad \bar{\Psi}_p \gamma_5 \Psi_n \pi^+$$

to be invariant under a global $U(1)$ transformation. That these terms conserve baryon number is evident since they contain only bilinear forms in the nucleon field. That they are also charge conserving follows from their invariance under the $U(1)_Q$ transformations

$$\begin{array}{rcl} \Psi_p \to \Psi_p' = \mathrm{e}^{-\mathrm{i}\epsilon q} \Psi_p & ; & \bar{\Psi}_p \to \bar{\Psi}_p' = \mathrm{e}^{+\mathrm{i}\epsilon q} \bar{\Psi}_p \\ \Psi_n \to \Psi_n' = \Psi_n & ; & \bar{\Psi}_n \to \bar{\Psi}_n' = \bar{\Psi}_n \\ \pi^+ \to (\pi^+)' = \mathrm{e}^{-\mathrm{i}\epsilon q} \pi^+ & ; & \pi^- \to (\pi^-)' = \mathrm{e}^{+\mathrm{i}\epsilon q} \pi^- \quad . \end{array}$$

The complete interaction Lagrangian now reads

$$\begin{aligned} \mathcal{L}_{\pi N} \;=\; & \mathrm{i}g_1 \bar{\Psi}_n \gamma_5 \Psi_n \pi^0 + \mathrm{i}g_2 \bar{\Psi}_p \gamma_5 \Psi_p \pi^0 \\ & + \mathrm{i}g_3 \bar{\Psi}_n \gamma_5 \Psi_p \pi^- + \mathrm{i}g_4 \bar{\Psi}_p \gamma_5 \Psi_n \pi^+ \quad . \end{aligned} \tag{3.70}$$

[5] Also possible is a coupling in which the derivative of the pion field appears ($\hat{Q} = \partial_\mu$). Since this is a pseudovector, it has to be multiplied by another pseudovector:

$$\mathcal{L}_{\pi N} \sim \bar{\Psi} \gamma_\mu \gamma_5 \Psi \, \partial^\mu \pi^0 \qquad \text{(pseudovector coupling)} \quad .$$

By introducing the matrices

$$\begin{aligned}
\frac{\tau_+}{2} &= \frac{1}{2}(\tau_1 + \mathrm{i}\tau_2) = \begin{pmatrix} 0 & 1 \\ 0 & 0 \end{pmatrix} \\
\frac{\tau_-}{2} &= \frac{1}{2}(\tau_1 - \mathrm{i}\tau_2) = \begin{pmatrix} 0 & 0 \\ 1 & 0 \end{pmatrix} \\
\tau_0 &= \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} = \tau_3
\end{aligned} \tag{3.71}$$

$\mathcal{L}_{\pi N}$ can be written in terms of the nucleon spinor Ψ as

$$\begin{aligned}
\mathcal{L}_{\pi N} &= \mathrm{i}g_1\bar{\Psi}\gamma_5\frac{1}{2}\left(1-\tau_0\right)\Psi\pi^0 + \mathrm{i}g_2\bar{\Psi}\gamma_5\frac{1}{2}\left(1+\tau_0\right)\Psi\pi^0 \\
&+\mathrm{i}g_3\bar{\Psi}\gamma_5\frac{\tau_-}{2}\Psi\pi^- + \mathrm{i}g_4\bar{\Psi}\gamma_5\frac{\tau_+}{2}\Psi\pi^+ \quad .
\end{aligned} \tag{3.72}$$

$\mathcal{L}_{\pi N}$ is, by construction, symmetric under $U(1)_B \otimes U(1)_Q$.

If now, in addition to charge and baryon number conservation, $SU(2)$ symmetry is also required for $\mathcal{L}_{\pi N}$, as suggested by the experimental fact of isospin conservation in the pion-nucleon interaction, then $\mathcal{L}_{\pi N}$ must be the scalar product of the isovector of the pion field with an isovector of the nucleon. Thus $\mathcal{L}_{\pi N}$ must have the structure

$$\mathcal{L}_{\pi N} = \mathrm{i}g\bar{\Psi}\gamma_5\frac{\boldsymbol{\tau}}{2}\Psi\cdot\boldsymbol{\pi} \quad , \tag{3.73}$$

with only one coupling constant, g. Since

$$\boldsymbol{\tau}\cdot\boldsymbol{\pi} = \frac{1}{\sqrt{2}}(\tau_+\pi^+ + \tau_-\pi^-) + \tau_0\pi^0 \quad ,$$

the Lagrangian of eq. (3.72) goes over into the form (3.73) for

$$g = -2g_1 = 2g_2 = \sqrt{2}g_3 = \sqrt{2}g_4 \quad .$$

There is then only one universal constant, g, that couples all pion fields to all states of the nucleon. This has the important consequence that the transition rates for seemingly quite different processes, like, e.g. , $p+p \to p+n+\pi^+$ and $n+n \to n+n+\pi^0$, are governed by one and the same coupling constant so that, if one rate has been measured, the other one can be predicted.

In summary, the interaction Lagrangian (3.73) is by construction invariant under baryon number and charge $U(1)$, isospin $SU(2)$, parity and

Lorentz transformations. In addition, it is also invariant under charge conjugation, i.e. particle-antiparticle conjugation. Its structure is that of a conserved current (cf. eq. (3.38)) multiplied with the pion field.

We have now the important result that the full pion-nucleon Lagrangian

$$\mathcal{L} = \mathcal{L}_N + \mathcal{L}_\pi + \mathcal{L}_{\pi N} \tag{3.74}$$

is $SU(2)$-invariant. As a consequence, not the isospin of nucleon and pion separately but instead the total isospin

$$\boldsymbol{T} = \boldsymbol{T}_N + \boldsymbol{T}_\pi = \int \Psi^\dagger \frac{\boldsymbol{\tau}}{2} \Psi \, d^3x - \mathrm{i} \int \partial_t \pi_k \, (\boldsymbol{t})_{kl} \, \pi_l \, d^3x \tag{3.75}$$

is conserved. The pion-nucleon coupling (3.73) does not contribute to $\boldsymbol{T}$ because it contains no field derivatives. Since $\mathcal{L}$ is obviously also $U(1)$ invariant, leading to baryon number conservation, it is also invariant under the combined $U(1) \otimes SU(2)$ transformation. Thus relation (3.45) linking charge, baryon-number and third component of isospin is also valid for the coupled system.

There is strong experimental evidence that $SU(2)$ is a good symmetry of the strong interactions so that the total isospin of any system of only strongly interacting particles is conserved [PER87]. If one takes into account additional fields that carry isospin, the expression (3.75) for $\boldsymbol{T}$ would contain contributions also from these other fields.

3.2.3 Structure of $SU(2)$ Multiplets

As already discussed in section 3.2.1, the basis states on which the isospin acts can be identified with particle multiplets. Since the algebra of $SU(2)$ is exactly that of the normal angular momentum in three-dimensional space it is evident that these multiplets can be characterized by the eigenvalues of $\boldsymbol{T}^2$; their individual states can be labeled by their eigenvalue of T_3. The mathematical reason for this is that $\boldsymbol{T}^2$ is the Casimir operator of the Lie group $SU(2)$. Since $SU(2)$ has only one Casimir operator, $\boldsymbol{T}^2$, it follows that besides $\boldsymbol{T}^2$ additionally only one generator, usually T_3, can be diagonalized simultaneously.

Starting from one state in a multiplet the others can be constructed by using the raising and lowering operators $T_+ = T_1 + \mathrm{i}T_2$ and $T_- = T_1 - \mathrm{i}T_2$. All the states within one multiplet have the same energy eigenvalue of the Hamiltonian, i.e. the same mass, if the underlying interaction is $SU(2)$-invariant. The $SU(2)$ transformation (3.56) transforms the states into others within the same multiplet of energetically degenerate states.

The product of two basis vectors of $SU(2)$ belonging to two multiplets (not necessarily different) does not in general form a basis vector of a specific multiplet but rather a superposition of basis vectors of other multiplets. For example, in the case of the pion-nucleon system treated in section 3.2.2 the nucleon forms an isospin doublet ($T = 1/2$) whereas the pions belong to an isospin triplet ($T = 1$). A simple nucleon-pion product state is then a superposition of states belonging to $T = 3/2$ and $T = 1/2$ multiplets. Their weight is given by the standard Clebsch-Gordan coefficients, well known from angular momentum coupling

$$|N\pi >_{T_3} = (|N > \otimes |\pi >)_{T_3} = \sum_T \left(\frac{1}{2} 1 T | T_{3N} T_{3\pi} T_3\right) |T T_3 > \quad , \tag{3.76}$$

where $T_3 = T_{3N} + T_{3\pi}$ and the sum over T runs over the values $|T| \geq |T_3|$. Thus the product of a triplet and doublet can be reduced into the sum of a doublet and a quadruplet, symbolically written as

$$3 \otimes 2 = 4 \oplus 2 \quad .$$

This reduction of the product can also be performed graphically. In this method all states of the product are represented by points on the T_3 axis. For each state of the product representation we have

$$T_3 = T_{3,1} + T_{3,2} \quad , \tag{3.77}$$

so that a given point T_3 can in general be constructed in several different ways. The highest possible value of T_3, however, is unique. Applying a lowering operator to this highest state generates the next lower state, and so on.

In the pion-nucleon example discussed above the highest state is the one with $T_3 = 3/2$, and the next lower is the one with $T_3 = 1/2$, which can be constructed in two different ways $((T_{3\pi} = 1) + (T_{3N} = -1/2)$ or $(T_{3\pi} = 0) + (T_{3N} = 1/2))$. One of these two states must belong to the $T = 3/2$ multiplet, the other is then assigned to the $T = 1/2$ multiplet. Graphically this looks as shown in Fig. 3.1.

The near equality of the masses within a given multiplet is a convincing evidence for the invariance of the strong interaction under $SU(2)$ transformations. The remaining mass splittings are, except for the pion, of the order of a few promille and can be ascribed to the electromagnetic interaction. The latter is true also for the pion; in this case the relative mass splitting is larger simply because of the exceptionally small mass of the pion.